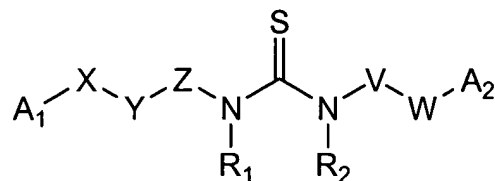


WHAT IS CLAIMED IS:

1. A compound of Formula 1



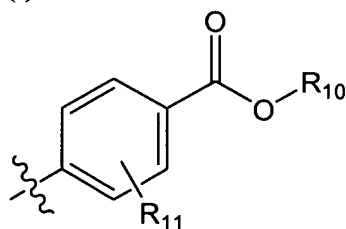
Formula 1

or a pharmaceutically acceptable salt thereof, wherein

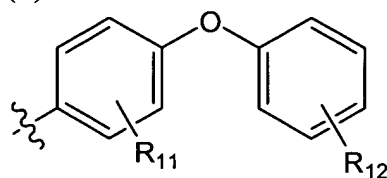
A₁ is an optionally substituted di-alkylamino, an optionally substituted aryl group, an optionally substituted 5- or 6- membered heteroaryl group, an optionally substituted bicyclic heteroaryl group having a 5-membered heteroaryl ring fused to a phenyl ring, an optionally substituted partially unsaturated or aromatic heterocyclic group having two 6-membered rings, an optionally substituted 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, an optionally substituted partially unsaturated 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, a 5- or 6-membered heterocycloalkyl group fused to a phenyl or heteroaryl ring, or a fused or spiro 8 to 11-membered bicyclic heterocycloalkyl group containing at least one nitrogen atom and 0 to 3 additional heteroatoms;

A₂ is

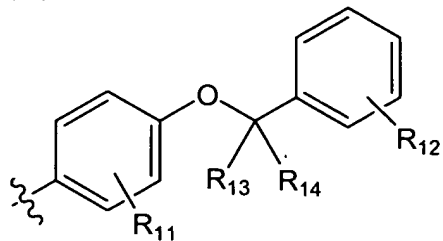
(i)



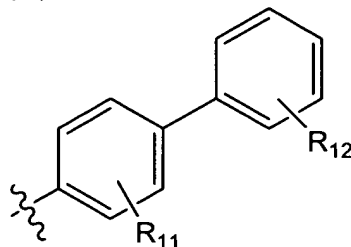
(ii)



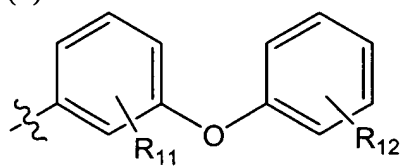
(iii)



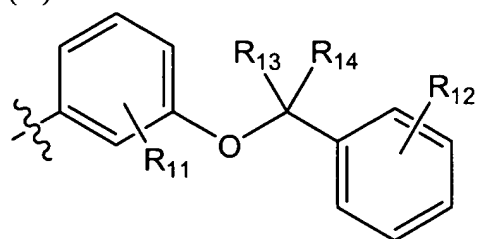
(iv)



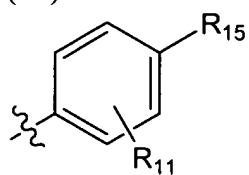
(v)



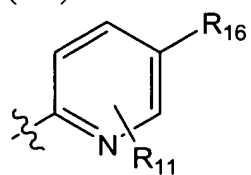
(vi)



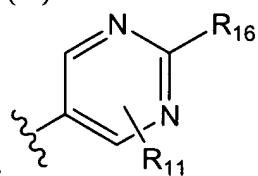
(vii)



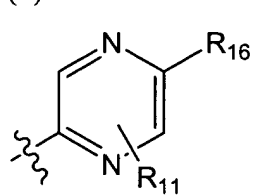
(viii)



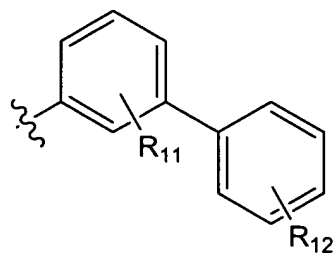
(ix)



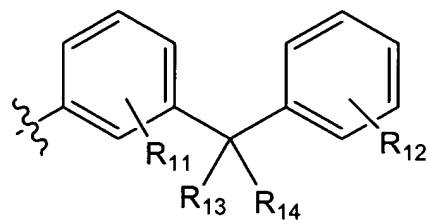
(x)



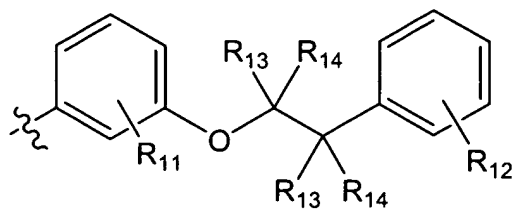
(xi)



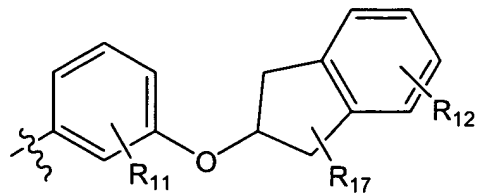
(xii)



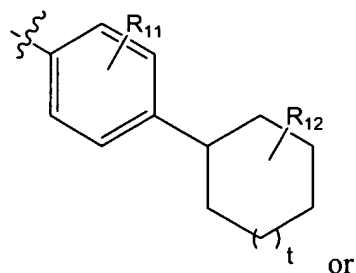
(xiii)



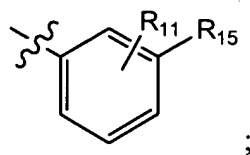
(xiv)



(xv)



(xvi)



t is 0 or 1;

X and W are independently O, S, NR, or absent, where R is hydrogen, optionally substituted C₁-C₆alkyl, or optionally substituted aryl(C₀-C₄alkyl);

V is C₁-C₆ alkyl, C₂-C₆alkenyl, C₃-C₇cycloalkyl, or absent;

Y is C₁-C₆ alkyl, C₁-C₆ alkyl substituted with C₃-C₇cycloalkyl, C₂-C₆alkenyl, C₃-C₇cycloalkyl, or absent;

wherein when V is absent, W is absent;

Z is carbonyl, thiocarbonyl, imino, or C₁-C₆alkylimino;

R₁ and R₂ are independently hydrogen, or

R₁ and R₂ are independently C₁-C₆alkyl, C₂-C₆ alkenyl, or C₂-C₆ alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkoxy, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy, or

R₁ and R₂ are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy;

R₁₀ is C₁-C₆alkyl;

R₁₁ and R₁₂ each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C₁-C₆alkyl, C₁-C₆alkoxy, mono- and di-(C₁-C₆alkyl)amino, C₂-C₆alkanoyl, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, and phenyl; and

R₁₃ and R₁₄ are independently chosen at each occurrence from hydrogen and C₁-C₄alkyl;

R₁₅ is C₄-C₆alkoxy or C₄-C₆alkyl;

R₁₆ is C₂-C₆alkoxy or C₂-C₆alkyl; and

R₁₇ represents 0 to 2 substituents independently chosen from halogen, methyl, and methoxy;

2. A compound or salt according to Claim 1 wherein

A₁ is a di-(C₁-C₆alkyl)amino, an aryl group, a 5- or 6- membered heteroaryl group, a bicyclic heteroaryl group having a 5-membered heteroaryl ring fused to a phenyl ring, a partially unsaturated or aromatic heterocyclic group having two 6-membered rings, a 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, a partially unsaturated 5- to 7-membered heterocycloalkyl group containing at least one nitrogen atom and 0 or 1 additional heteroatoms, a 5- or 6-membered heterocycloalkyl group fused to a phenyl or heteroaryl ring, or a fused or spiro 8 to 11-membered bicyclic heterocycloalkyl group containing at least one nitrogen atom and 0 to 3 additional heteroatoms; each of which A₁ is substituted with 0 to 5 substituents independently chosen from:

- (a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -CONH₂, -SO₂NH₂, -SH, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy,
- (b) C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₁-C₆alkoxy, C₂-C₆alkenyloxy, C₁-C₄alkoxy(C₁-C₄alkyl), amino(C₁-C₆alkyl), mono- and di-(C₁-C₆alkyl)amino, mono- and di-(C₁-C₄alkyl)aminoC₁-C₄alkyl, C₂-C₆alkanoyl, C₂-C₈alkanoyloxy, C₁-C₈alkoxycarbonyl, -mono- and di-(C₁-C₆alkyl)carboxamide, (C₃-C₇cycloalkyl)carboxamide, mono- and di-(C₁-C₆alkyl)sulfonamide, C₁-C₆alkylthio, aryl(C₀-C₄alkyl)thio, C₁-C₆alkylsulfinyl, and C₁-C₆alkylsulfonyl, and
- (c) -GR_a where

G is chosen from -(CH₂)_n-, C₂-C₄alkenyl, C₂-C₄alkynyl, -(CH₂)_nO(CH₂)_m-, and -(CH₂)_nN(CH₂)_m-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C₃-C₈cycloalkyl, C₂-C₇monocyclic heterocycloalkyl, C₅-C₁₀bicyclic heterocycloalkyl, indanyl, tetrahydronaphthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₂-C₄alkanoyl, C₁-C₄alkoxycarbonyl, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, and phenyl;

X and W are independently O, S, NR, or absent,

where R is hydrogen or R is C₁-C₆alkyl or aryl(C₀-C₄alkyl), each of which is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, C₁-C₆alkyl, C₁-C₆alkoxy, and mono- and di-(C₁-C₆alkyl)amino;

V is independently C₁-C₆ alkyl, C₂-C₆alkenyl, C₃-C₇cycloalkyl, or absent;

Y is C₁-C₆ alkyl, C₁-C₆ alkyl substituted with C₃-C₇cycloalkyl, C₂-C₆alkenyl, C₃-C₇cycloalkyl, or absent;

wherein when V is absent, W is absent;

Z is carbonyl, thiocarbonyl, or imino; and

R₁ and R₂ are independently hydrogen, or

R₁ and R₂ are independently C₁-C₆alkyl, C₂-C₆ alkenyl, or C₂-C₆ alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkoxy, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy, or

R₁ and R₂ are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy.

3. A compound or salt according to Claim 1 or 2 in which Z is thiocarbonyl.
4. A compound or salt according to Claim 1 or 2 in which Z is imino or C₁-C₆alkylimino.
5. A compound or salt according to Claim 4 in which Z is imino or methylimino.
6. A compound or salt according to Claim 1 or 2 in which Z is carbonyl.
7. A compound or salt according to any one of Claims 1 to 6 in which X is oxygen and Y is -CH₂-.

8. A compound or salt according to any one of Claims 1 to 6 in which X is oxygen and Y is $-\text{CH}_2\text{CH}_2-$.
9. A compound or salt according to any one of Claims 1 to 6 wherein when X and Y are absent.
10. A compound or salt according to any one of Claims 1 to 9 wherein V and W are absent.
11. A compound or salt according to any one of Claims 1 to 9 in which V is C_1 - C_2 alkyl and W is absent.
12. A compound or salt according to any one of Claims 1 to 11 in which R_1 and R_2 are independently hydrogen, or C_1 - C_4 alkyl, C_2 - C_4 alkenyl, or C_2 - C_4 alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1 - C_4 alkoxy, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.
13. A compound or salt according to Claim 12 in which R_1 and R_2 are independently hydrogen, methyl, or ethyl.
14. A compound or salt according to Claim 13 in which R_1 and R_2 are both hydrogen.
15. A compound or salt according to any one of Claims 1 to 11 in which R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(\text{C}_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

16. A compound or salt according to Claim 15 in which R_1 and R_2 are joined to form a 5- to 7-membered saturated or mono-unsaturated ring containing no additional heteroatoms, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C_1 - C_2 alkyl, and C_1 - C_2 alkoxy.

17. A compound or salt according to any one of Claims 2 to 16 wherein A_1 is aryl, a partially unsaturated heterocyclic group, or heteroaryl group; substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, $-COOH$, $-CONH_2$, $-SO_2NH_2$, $-SH$, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy, and

(b) C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_1 - C_6 alkoxy, C_2 - C_6 alkenyloxy, C_1 - C_4 alkoxy(C_1 - C_4 alkyl), amino(C_1 - C_6 alkyl), mono- and di-(C_1 - C_6 alkyl)amino, mono- and di-(C_1 - C_4 alkyl)amino, C_1 - C_4 alkyl, C_2 - C_6 alkanoyl, C_2 - C_8 alkanoyloxy, C_1 - C_8 alkoxycarbonyl, mono- and di-(C_1 - C_6 alkyl)carboxamide, (C_3 - C_7 cycloalkyl)carboxamide, mono- and di-(C_1 - C_6 alkyl)sulfonamide, C_1 - C_6 alkylthio, aryl(C_0 - C_4 alkyl)thio, C_1 - C_6 alkylsulfinyl, and C_1 - C_6 alkylsulfonyl, and

(c) $-GR_a$ where

G is chosen from $-(CH_2)_n-$, C_2 - C_4 alkenyl, C_2 - C_4 alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3 - C_8 cycloalkyl, C_2 - C_7 monocyclic heterocycloalkyl, C_5 - C_{10} bicyclic heterocycloalkyl, indanyl, tetrahydronaphthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di-(C_1 - C_4 alkyl)amino, C_2 - C_4 alkanoyl, C_1 - C_4 alkoxycarbonyl, C_1 - C_2 haloalkyl, C_1 - C_2 haloalkoxy, and phenyl.

18. A compound or salt according to Claim 17 wherein

A₁ is phenyl, naphthyl, pyridyl, pyrimidinyl, thienyl, pyrrolyl, furanyl, pyrazolyl, imidazolyl, thiazolyl, triazolyl, thiadiazolyl, oxazolyl, isoxazolyl, benzofuranyl, benzothiazolyl, benzothiophenyl, benzoxadiazolyl, benzo[*d*]oxazolyl, dihydrobenzodioxynyl, indolyl, pyrazolopyrimidinyl, thienylpyrazolyl, benzopyranyl, or 4*H*-chromenyl,

each of which is substituted with 0 to 5 substituents independently chosen from

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -CONH₂, -SO₂NH₂, -SH, C₁-

C₂haloalkyl, and C₁-C₂haloalkoxy, and

(b) C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₁-C₆alkoxy, C₂-C₆alkenyloxy, C₁-C₄alkoxy(C₁-C₄alkyl), amino(C₁-C₆)alkyl, mono- and di-(C₁-C₆alkyl)amino, mono- and di-(C₁-C₄alkyl)aminoC₁-C₄alkyl, C₂-C₆alkanoyl, C₂-C₈alkanoyloxy, C₁-C₈alkoxycarbonyl, -mono- and di-(C₁-C₆alkyl)carboxamide, (C₃-C₇cycloalkyl)carboxamide, mono- and di-(C₁-C₆alkyl)sulfonamide, C₁-C₆alkylthio, aryl(C₀-C₄alkyl)thio, C₁-C₆alkylsulfinyl, and C₁-C₆alkylsulfonyl, and

(c) -GR_a where

G is chosen from -(CH₂)_n-, C₂-C₄alkenyl, C₂-C₄alkynyl, -(CH₂)_nO(CH₂)_m-, and -(CH₂)_nN(CH₂)_m-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C₃-C₈cycloalkyl, C₂-C₇monocyclic heterocycloalkyl, C₅-

C₁₀bicyclicheterocycloalkyl, indanyl, tetrahydronaphthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from

halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, and phenyl.

19. A compound or salt according to Claim 18 in which

A₁ is phenyl, naphthyl, pyridin-2-yl, pyridin-3-yl, pyridin-4-yl, pyrimidin-2-yl, pyrimidin-4-yl, pyrimidin-5-yl, thien-2-yl, thien-3-yl, thiazol-4-yl, pyrrol-1-yl, pyrrol-2-yl, pyrrol-3-yl, furan-2-yl, furan-3-yl, pyrazol-1-yl, pyrazol-2-yl, pyrazol-4-yl, pyrazol-5-yl, imidazol-1-yl, imidazol-2-yl, imidazol-4-yl, imidazol-5-yl, thiazol-2-yl, thiazol-3-yl, thiazol-5-yl, 1,2,3-triazol-4-yl, 1,2,3-thiadiazol-4-yl, 1,2,3-thiadiazol-5-yl, oxazol-2-yl, isoxazol-4-yl, isoxazol-5-yl, isoxazol-3-yl, isoxazol-4-yl, isoxazol-5-yl, benzofuran-2-yl, benzofuran-3-yl, benzopyran-2-yl, benzopyran-3-yl, benzopyran-4-yl, benzo[*d*]oxazol-2-yl, benzo[*d*]thiazol-2-yl, benzo[*b*]thiophen-2-yl, 4*H*-chromen-2-yl, benzo[*c*][1,2,5]oxadiazolyl, 2,3-dihydrobenzo[*b*][1,4]dioxin-2-yl, pyrazolo[1,5-*a*]pyrimidin-6-yl, dihydrobenzo[*b*][1,4]dioxin-3-yl, indol-2-yl, pyrazolo[1,5-*a*]pyrimidin-5-yl, 1*H*-thieno[2,3-*c*]pyrazol-4-yl, or 1*H*-thieno[2,3-*c*]pyrazol-5-yl,

each of which is substituted with 0 to 5 substituents independently chosen from

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -CONH₂, -SO₂NH₂, -SH, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy, and

(b) C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₁-C₆alkoxy, C₂-C₆alkenyloxy, C₁-C₄alkoxy(C₁-C₄alkyl), amino(C₁-C₆alkyl), mono- and di-(C₁-C₆alkyl)amino, mono- and di-(C₁-C₄alkyl)amino, C₁-C₄alkyl, C₂-C₆alkanoyl, C₂-C₈alkanoyloxy, C₁-C₈alkoxycarbonyl, -mono- and di-(C₁-C₆alkyl)carboxamide, (C₃-C₇cycloalkyl)carboxamide, mono- and di-(C₁-C₆alkyl)sulfonamide, C₁-C₆alkylthio, aryl(C₀-C₄alkyl)thio, C₁-C₆alkylsulfinyl, and C₁-C₆alkylsulfonyl, and

(c) -GR_a where

G is chosen from -(CH₂)_n-, C₂-C₄alkenyl, C₂-C₄alkynyl, -(CH₂)_nO(CH₂)_m-, and -(CH₂)_nN(CH₂)_m-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C₃-C₈cycloalkyl, C₂-C₇monocyclic heterocycloalkyl, C₅-

C₁₀bicyclic heterocycloalkyl, indanyl, tetrahydronaphthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from

halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, and phenyl.

20. A compound or salt according to Claim 19 in which A₁ is substituted with 0 to 5 substituents independently chosen from

- (a) halogen, hydroxy, cyano, amino, nitro, oxo, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy, and
- (b) C₁-C₆alkyl, C₁-C₆alkoxy, C₁-C₄alkoxy(C₁-C₄alkyl), amino(C₁-C₄)alkyl, mono- and di-(C₁-C₄alkyl)amino, and mono- and di-(C₁-C₄alkyl)aminoC₁-C₄alkyl; and
- (c) -GR_a where

G is from -(CH₂)_n-, -(CH₂)_nO(CH₂)_m-, and -(CH₂)_nN(CH₂)_m-, and

R_a is C₃-C₈cycloalkyl, 5 or 6-membered heterocycloalkyl containing 1 or 2 heteroatoms independently chosen from O, S, and N, 5- or 6-membered heteroaryl containing 1, 2, or 3 heteroatoms independently chosen from O, S, and N, indanyl, and phenyl,

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₂alkyl)amino, and C₁-C₂haloalkyl, and C₁-C₂haloalkoxy.

21. A compound or salt according to Claim 20 in which

A₁ is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, C₁-C₆alkyl, C₁-C₆alkoxy, C₁-C₄alkoxy(C₁-C₄alkyl), amino(C₁-C₄)alkyl, mono- and di-(C₁-C₄alkyl)amino, and mono- and di-(C₁-C₄alkyl)aminoC₁-C₄alkyl.

22. A compound or salt according to any one of Claim 3 to 16 wherein A_1 is C_1 - C_6 alkyl, C_3 - C_7 cycloalkyl, or C_2 - C_7 monocyclic heterocycloalkyl, each of which is substituted with 0 to 5 substituents independently chosen from:
- (a) halogen, hydroxy, cyano, amino, nitro, oxo, $-COOH$, $-CONH_2$, $-SO_2NH_2$, $-SH$, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy, and
- (b) C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_1 - C_6 alkoxy, C_2 - C_6 alkenyloxy, C_1 - C_4 alkoxy(C_1 - C_4 alkyl), amino(C_1 - C_6)alkyl, mono- and di-(C_1 - C_6 alkyl)amino, mono- and di-(C_1 - C_4 alkyl)amino C_1 - C_4 alkyl, C_2 - C_6 alkanoyl, C_2 - C_8 alkanoyloxy, C_1 - C_8 alkoxycarbonyl, -mono- and di-(C_1 - C_6 alkyl)carboxamide, (C_3 - C_7 cycloalkyl)carboxamide, mono- and di-(C_1 - C_6 alkyl)sulfonamide, C_1 - C_6 alkylthio, aryl(C_0 - C_4 alkyl)thio, C_1 - C_6 alkylsulfinyl, and C_1 - C_6 alkylsulfonyl, and
- (c) $-GR_a$ where
- G is chosen from $-(CH_2)_n-$, C_2 - C_4 alkenyl, C_2 - C_4 alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and
- R_a is chosen from C_3 - C_8 cycloalkyl, C_2 - C_7 monocyclic heterocycloalkyl, C_5 - C_{10} bicyclic heterocycloalkyl, indanyl, tetrahydronaphthyl, aryl, and heteroaryl;
- each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di-(C_1 - C_4 alkyl)amino, C_2 - C_4 alkanoyl, C_1 - C_4 alkoxycarbonyl, C_1 - C_2 haloalkyl, C_1 - C_2 haloalkoxy, and phenyl.

23. A compound or salt according to Claim 22 wherein

A₁ is C₁-C₆ alkyl, C₃-C₇cycloalkyl, pyrrolidinyl, piperidinyl, piperazinyl, or morpholinyl; each of which is substituted with 0 to 3 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -CONH₂, -SO₂NH₂, -SH, C₁-

C₂haloalkyl, and C₁-C₂haloalkoxy, and

(b) C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₁-C₆alkoxy, C₂-C₆alkenyloxy, C₁-C₄alkoxy(C₁-C₄alkyl), amino(C₁-C₆alkyl), mono- and di-(C₁-C₆alkyl)amino, mono- and di-(C₁-C₄alkyl)aminoC₁-C₄alkyl, C₂-C₆alkanoyl, C₂-C₈alkanoyloxy, C₁-C₈alkoxycarbonyl, -mono- and di-(C₁-C₆alkyl)carboxamide, (C₃-C₇cycloalkyl)carboxamide, mono- and di-(C₁-C₆alkyl)sulfonamide, C₁-C₆alkylthio, aryl(C₀-C₄alkyl)thio, C₁-C₆alkylsulfinyl, and C₁-C₆alkylsulfonyl, and

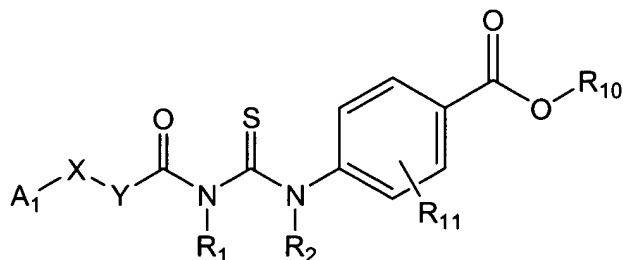
(c) -GR_a where

G is chosen from -(CH₂)_n-, C₂-C₄alkenyl, C₂-C₄alkynyl, -(CH₂)_nO(CH₂)_m-, and

-(CH₂)_nN(CH₂)_m-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C₃-C₈cycloalkyl, and phenyl.

24. A compound or salt according to Claim 1 or 2 of Formula 2, wherein :



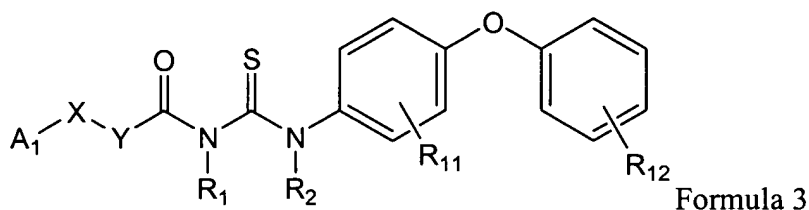
Formula 2

R₁ and R₂ are hydrogen or methyl;

R₁₀ is C₁-C₆alkyl; and

R₁₁ represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy.

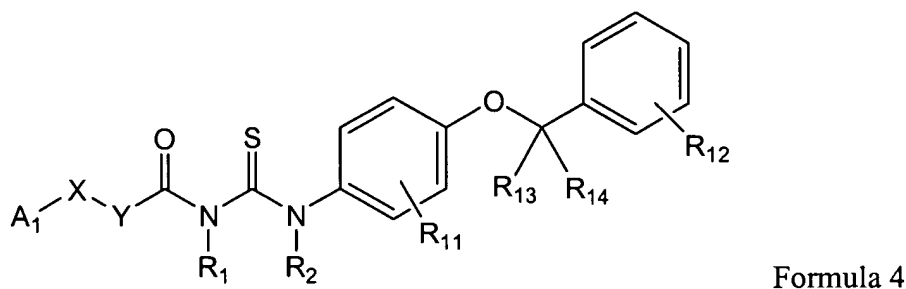
25. A compound or salt according to Claim 1 or 2 of Formula 3, wherein:



R_1 and R_2 are hydrogen or methyl; and

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

26. A compound or salt according to Claim 1 or 2 of Formula 4, wherein:

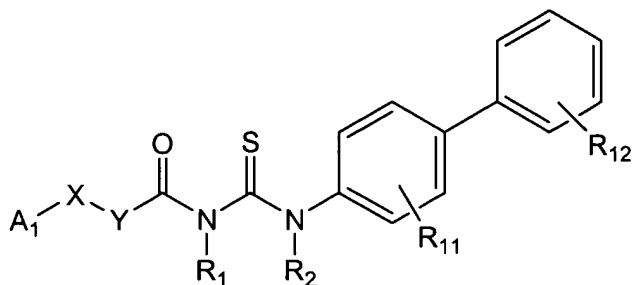


R_1 and R_2 are independently hydrogen or methyl;

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

R_{13} and R_{14} are independently hydrogen or methyl.

27. A compound or salt according to Claim 1 or 2 of Formula 5, wherein

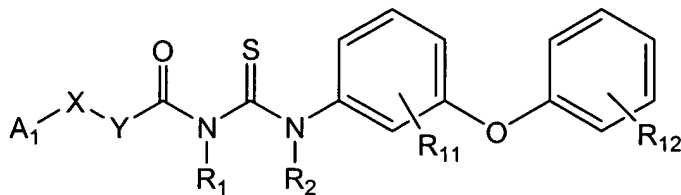


Formula 5

R_1 and R_2 are independently hydrogen or methyl; and

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

28. A compound or salt according to any one of Claim 1 or 2 of Formula 6, wherein

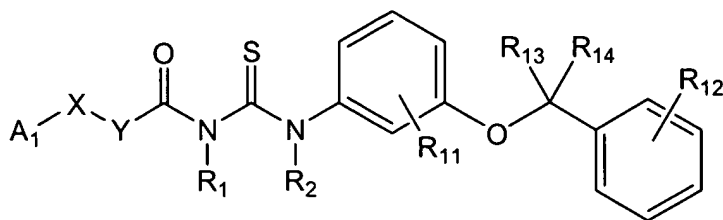


Formula 6

R_1 and R_2 are independently hydrogen or methyl; and

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

29. A compound or salt according to Claim 1 or 2 of Formula 7:



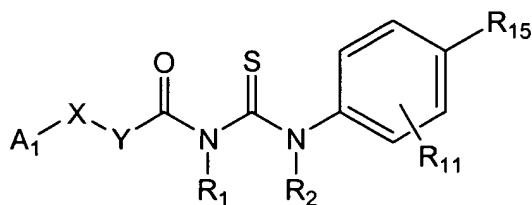
Formula 7

R₁ and R₂ are independently hydrogen or methyl;

R₁₁ and R₁₂ each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy; and

R₁₃ and R₁₄ are independently hydrogen or methyl.

30. A compound or salt according to Claim 1 or 2 of Formula 8, wherein



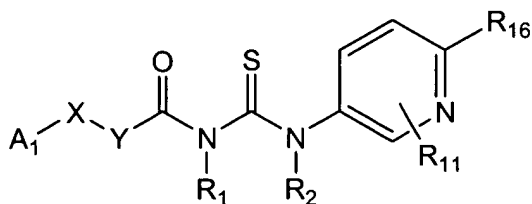
Formula 8

R₁ and R₂ are independently hydrogen or methyl;

R₁₁ represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy; and

R₁₅ represents C₄-C₆alkoxy or C₄-C₆alkyl.

31. A compound salt according to Claim 1 or 2 of Formula 9, wherein



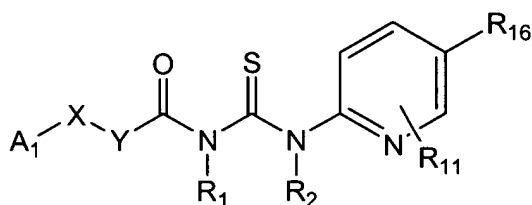
Formula 9

R_1 and R_2 are independently hydrogen or methyl;

R_{11} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

R_{16} is C_2 - C_6 alkoxy or C_2 - C_6 alkyl.

32. A compound or salt according to Claim 1 or 2 of Formula 10, wherein



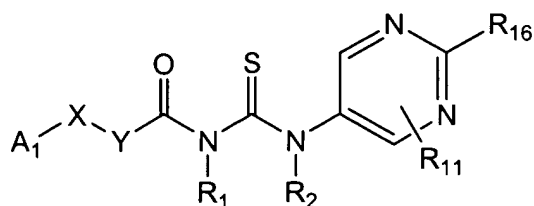
Formula 10

R_1 and R_2 are independently hydrogen or methyl;

R_{11} is represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

R_{16} is C_2 - C_6 alkoxy or C_2 - C_6 alkyl.

33. A compound or salt according to Claim 1 or 2 of Formula 11, wherein



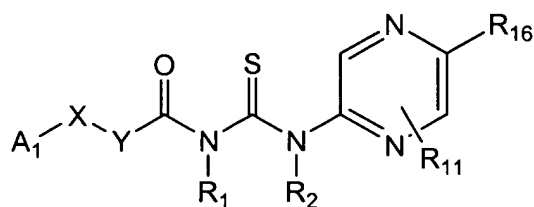
Formula 11

R_1 and R_2 are independently hydrogen or methyl;

R_{11} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

R_{16} is C_2 - C_6 alkoxy or C_1 - C_6 alkyl.

34. A compound or salt according to Claim 1 or 2 of Formula 12, wherein



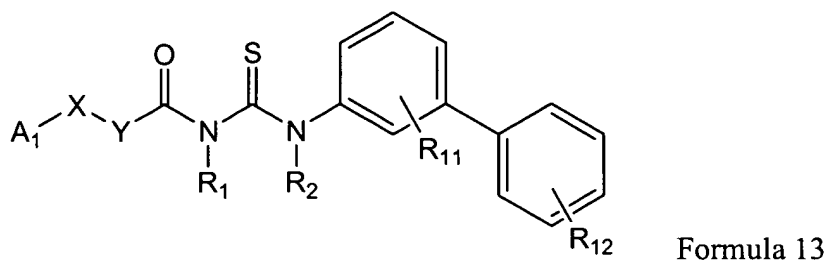
Formula 12

R_1 and R_2 are independently hydrogen or methyl;

R_{11} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

R_{16} is C_2 - C_6 alkoxy or C_2 - C_6 alkyl.

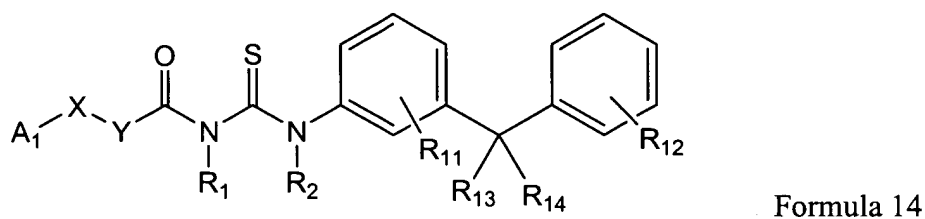
35. A compound or salt according to Claim 1 or 2 of Formula 13, wherein



R_1 and R_2 are independently hydrogen or methyl; and

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

36. A compound or salt according to Claim 1 or 2 of Formula 14, wherein

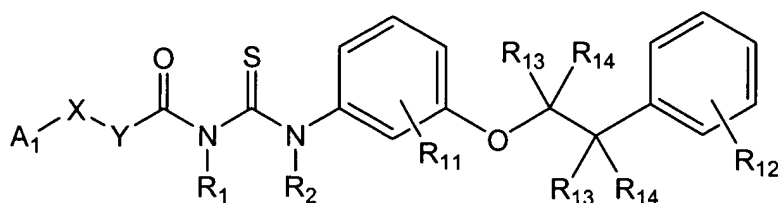


R_1 and R_2 are independently hydrogen or methyl;

R_{11} and R_{12} each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

R_{13} and R_{14} are independently hydrogen or methyl.

37. A compound salt according to Claim 1 or 2 of Formula 15, wherein



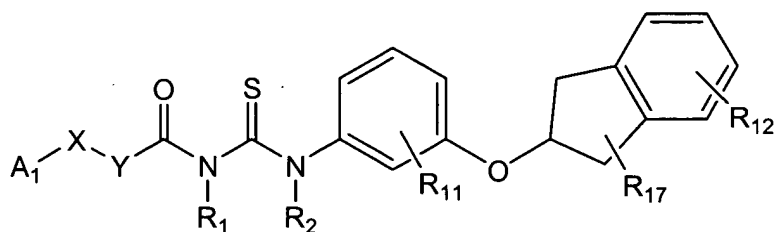
Formula 15

R₁ and R₂ are independently hydrogen or methyl;

R₁₁ and R₁₂ each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy; and

R₁₃ and R₁₄ are independently chosen at each occurrence from hydrogen and methyl.

38. A compound salt according to Claim 1 or 2 of Formula 16, wherein



Formula 16

R₁ and R₂ are independently hydrogen or methyl;

R₁₁ and R₁₂ each represent 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy; and

R₁₇ represents 0 to 2 substituents independently chosen from halogen, methyl, and methoxy.

39. A compound or salt according to any one of Claim 24 to 38 in which

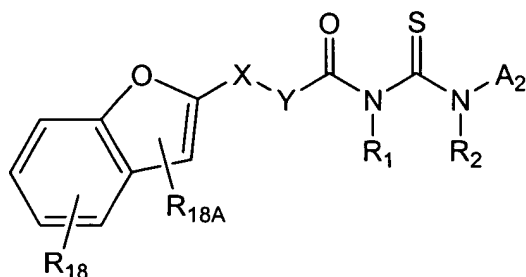
X is NR and Y is -CH₂- or -CH₂CH₂-; or

X is O and Y is -CH₂- or -CH₂CH₂-; or

X and Y are absent.

40. A compound or salt according to any one of Claims 1, 2, or 24 to 38, wherein A_1 is pyrazinyl, pyridyl, or quinaxolynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di-(C_1 - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

41. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 17, wherein



Formula 17

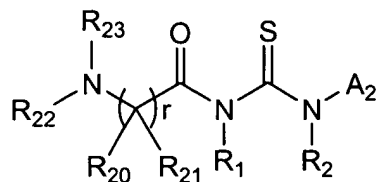
wherein

R_{18A} is hydrogen, halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di-(C_1 - C_4 alkyl)amino, C_1 - C_2 haloalkyl, or C_1 - C_2 haloalkoxy; and

R_{18} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di-(C_1 - C_4 alkyl)amino, C_1 - C_2 haloalkyl, C_1 - C_2 haloalkoxy, and phenyl.

42. A compound or salt according to Claim 41 in which X and Y are absent; and R_1 and R_2 are independently hydrogen or methyl.

43. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 18, wherein



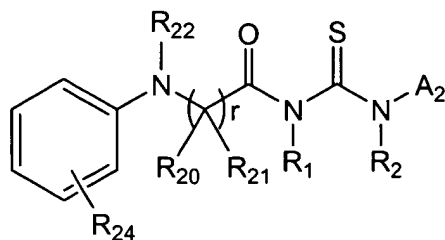
Formula 18

r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1 - C_4 alkyl; or R_{20} and R_{21} are joined to form a C_3 - C_7 cycloalkyl group; and

R_{22} and R_{23} are independently chosen C_1 - C_6 alkyl groups; each of which is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

44. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 19, wherein



Formula 19

r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1 - C_4 alkyl; or R_{20} and R_{21} are joined to form a C_3 - C_7 cycloalkyl group;

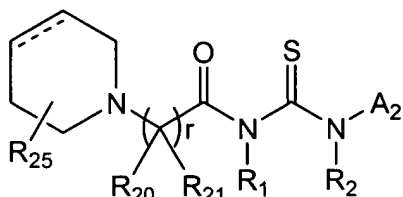
R_{22} is C_1 - C_6 alkyl which is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

R_{24} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, $-COOH$, $-CONH_2$, $-SO_2NH_2$, $-SH$, C_1 - C_2 haloalkyl, C_1 - C_2 haloalkoxy, C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_1 - C_6 alkoxy, C_2 - C_6 alkenyloxy, C_1 - C_4 alkoxy(C_1 - C_4 alkyl), amino(C_1 - C_6 alkyl), mono- and di- $(C_1$ - C_6 alkyl)amino, mono- and di- $(C_1$ - C_4 alkyl)amino C_1 - C_4 alkyl, C_2 - C_6 alkanoyl, C_2 - C_8 alkanoyloxy, C_1 - C_8 alkoxycarbonyl, -mono- and di- $(C_1$ - C_6 alkyl)carboxamide, $(C_3$ - C_7 cycloalkyl)carboxamide, mono- and di- $(C_1$ - C_6 alkyl)sulfonamide, C_1 - C_6 alkylthio, aryl(C_0 - C_4 alkyl)thio, C_1 - C_6 alkylsulfinyl, and C_1 - C_6 alkylsulfonyl.

45. A compound or salt according to Claim 44, wherein

R_{24} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

46. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 20, wherein



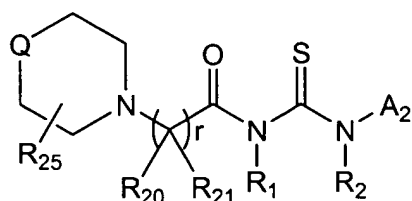
Formula 20

r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1 - C_4 alkyl; or R_{20} and R_{21} are joined to form a C_3 - C_7 cycloalkyl group; and

R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

47. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 21, wherein



Formula 21

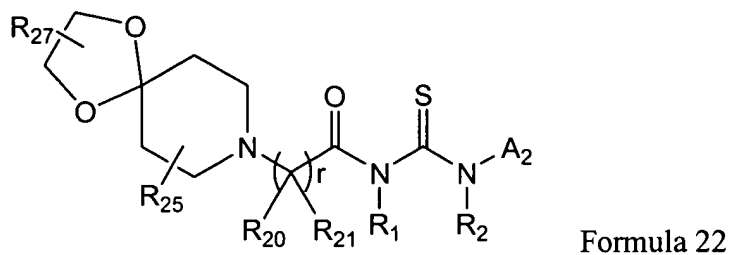
r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1 - C_4 alkyl; or R_{20} and R_{21} are joined to form a C_3 - C_7 cycloalkyl group; and

R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

Q is O, S, or NR_{26} ; where R_{26} is hydrogen or R_{26} is C_1 - C_6 alkyl, phenyl, pyridyl, or pyrimidinyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_6 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

48. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 22, wherein

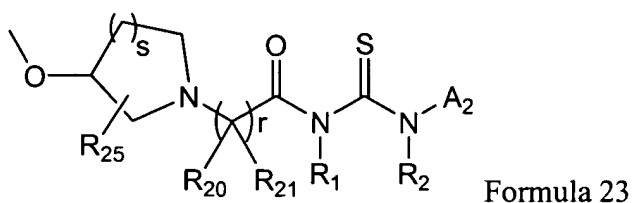


r is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1 - C_4 alkyl; or R_{20} and R_{21} are joined to form a C_3 - C_7 cycloalkyl group; and

R_{25} and R_{27} each represent 0 to 2 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

49. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 23, wherein



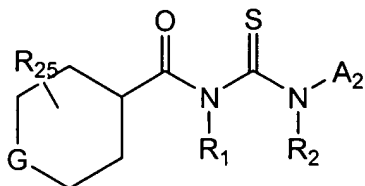
r is 1, 2, or 3;

s is 1, 2, or 3;

R_{20} and R_{21} are independently selected from hydrogen and C_1 - C_4 alkyl; or R_{20} and R_{21} are joined to form a C_3 - C_7 cycloalkyl group; and

R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

50. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 24, wherein



Formula 24

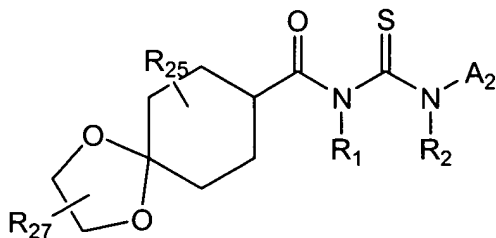
R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_6 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

G is O, S, SO_2 , or NR_{26} ; where R_{26} is hydrogen or R_{26} is C_1 - C_6 alkyl, phenyl, pyridyl, or pyrimidinyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_6 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

51. A compound or salt according to Claim 50, wherein

R_{25} represents a di- $(C_1$ - C_6 alkyl)amino substituent and 0 to 2 additional substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_6 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

52. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 25, wherein



Formula 25

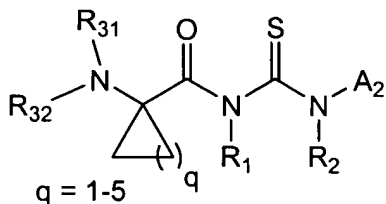
R_{25} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

R_{27} represents 0 to 2 substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

53. A compound or salt according to Claim 52, wherein

R_{25} represents a di- $(C_1$ - C_6 alkyl)amino substituent and 0 to 2 additional substituents independently chosen from halogen, hydroxy, cyano, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_6 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

54. A compound or salt according any one of Claims 1, 2 or 24 to 38 of Formula 26, wherein

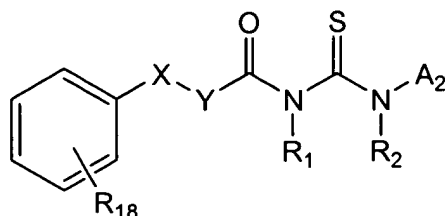


Formula 26

q is an integer from 1 to 5; and

R_{31} and R_{32} are independently chosen from C_1 - C_6 alkyl and phenyl; each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy.

55. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 27, wherein



Formula 27

wherein

R_{18} represents 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, C_1 - C_2 haloalkoxy, and phenyl.

56. A compound or salt according to Claim 55 wherein X is oxygen and Y is $-CH_2-$.

57. A compound or salt according to Claim 55 wherein X and Y are absent.

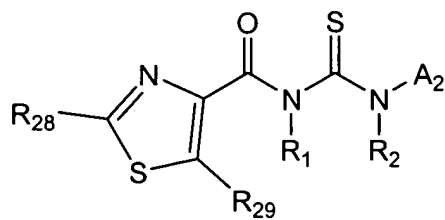
58. A compound or salt according to any one of Claims 1, 2, or 24 to 38, wherein A_1 is 5-membered heteroaryl group selected from furan-2-yl, furan-3-yl, isoxazol-3-yl, isoxazol-4-yl, thiophen-2-yl, thiophen-3-yl, pyrrol-2-yl, pyrrol-3-yl, and pyrazolyl; each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1 - C_4 alkyl, C_2 - C_4 alkenyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, C_1 - C_2 haloalkoxy, and phenyl.

59. A compound or salt according to Claim 58 wherein X is oxygen and Y is $-CH_2-$.

60. A compound or salt according to Claim 58 wherein X and Y are absent.

61. A compound or salt according to any one of Claims 1, 2, or 24 to 38, wherein A_1 is pyridin-2-yl or pyridin-3-yl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, C_1 - C_2 haloalkoxy, and phenyl.

62. A compound or salt according to Claim 61 wherein X is oxygen and Y is $-\text{CH}_2-$.
63. A compound or salt according to Claim 61 wherein X and Y are absent.
64. A compound or salt according to any one of Claims 1, 2, or 24 to 38 of Formula 28, wherein



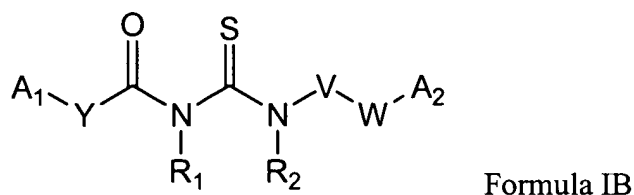
Formula 28

wherein

R_{28} is phenyl or pyridyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(\text{C}_1$ - C_4 alkyl)amino, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy; and

R_{29} is hydrogen, methyl or ethyl.

65. A compound or pharmaceutically acceptable salt of Formula IB, wherein



A₁ is di-(C₁-C₈alkyl)amino, an *N*-(C₁-C₆alkyl)-*N*-phenyl-amino group, an *N*-(C₁-C₆alkyl)-*N*-pyridyl amino group, a 5- to 7-membered monocyclic heterocycloalkyl group covalently bound to a point of attachment in Formula IB via a Nitrogen atom, a 5- to 7-membered monocyclic partially unsaturated heterocyclic group covalently bound to a point of attachment in Formula IB via a Nitrogen atom, a 5- to 7- membered heterocycloalkyl group covalently bound to a point of attachment in Formula IB via a Carbon atom which is adjacent to a Nitrogen atom, or an 8- to 11- membered bicyclic heterocycloalkyl in which the rings are fused or spiro covalently bound to a point of attachment in Formula IB via a Nitrogen atom;

A₂ is C₃-C₈ cycloalkyl, a partially unsaturated or aromatic carbocyclic group, or a saturated, partially unsaturated, or aromatic heterocyclic group;

each of which A₁ and A₂ is substituted with 0 to 5 substituents independently chosen from (a), (b), and (c), where

- (a) is independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -CONH₂, -SO₂NH₂, -SH, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy,
- (b) is independently chosen from C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₁-C₆alkoxy, C₂-C₆alkenyloxy, C₁-C₄alkoxy(C₁-C₄alkyl), amino(C₁-C₆alkyl), mono- and di-(C₁-C₆alkyl)amino, mono- and di-(C₁-C₄alkyl)aminoC₁-C₄alkyl, C₂-C₆alkanoyl, C₂-C₈alkanoyloxy, C₁-C₈alkoxycarbonyl, -mono- and di-(C₁-C₆alkyl)carboxamide, (C₃-C₇cycloalkyl)carboxamide, mono- and di-(C₁-C₆alkyl)sulfonamide, C₁-C₆alkylthio, aryl(C₀-C₄alkyl)thio, C₁-C₆alkylsulfinyl, and C₁-C₆alkylsulfonyl, and
- (c) is -GR_a where G is chosen from -(CH₂)_n-, C₂-C₄alkenyl, C₂-C₄alkynyl, -(CH₂)_nO(CH₂)_m-, and -(CH₂)_nN(CH₂)_m-, where n and m are independently 0, 1, 2, 3, or 4; and
- R_a is chosen from C₃-C₈cycloalkyl, C₂-C₇monocyclic heterocycloalkyl, C₅-C₁₀bicyclic heterocycloalkyl, indanyl, tetrahydronaphthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₂-C₄alkanoyl, C₁-C₄alkoxycarbonyl, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, and phenyl;

W is O, S, NR, or absent, where R is hydrogen or R is C₁-C₆alkyl or aryl(C₀-C₄alkyl), each of which is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, oxo, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, C₁-C₆alkyl, C₁-C₆alkoxy, and mono- and di-(C₁-C₆alkyl)amino;

V is C₁-C₆ alkyl, C₂-C₆alkenyl, C₃-C₇cycloalkyl, or absent; and when V is absent, W is absent;

Y is C₁-C₆ alkyl substituted with

0 or 1 of C₃-C₇cycloalkyl, a 5- to 7-membered monocyclic heterocycloalkyl, or 8- to 11-membered bicyclic heterocycloalkyl in which the rings are fused or spiro; each of which substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, cyano, amino, nitro, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy; or

Y is absent;

R₁ and R₂ are independently hydrogen or C₁-C₆alkyl, C₂-C₆ alkenyl, or C₂-C₆ alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkoxy, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy, or R₁ and R₂ are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy.

66. A compound or salt according to Claim 65 wherein V and W are absent.

67. A compound or salt according to Claim 65 in which Y is absent.

68. A compound or salt according to Claim 65 in which Y is -CH₂-.

69. A compound or salt according to Claim 65 in which Y is $-\text{CH}_2-$ substituted with C_3 - C_6 cycloalkyl, pyrrolidinyl, or piperidinyl.

70. A compound or salt according to any one of Claims 65 to 69 in which R_1 and R_2 are independently hydrogen or C_1 - C_4 alkyl.

71. A compound or salt according to Claim 70 in which R_1 and R_2 are independently hydrogen or methyl.

72. A compound or salt according to any one of Claims 65 to 71 wherein A_2 is C_5 - C_7 cycloalkyl, phenyl, pyridyl, naphthyl, pyrimidinyl, pyrazinyl, benzothiazolyl, benzodioxyl, quinolinyl, or isoquinolinyl, each of which is substituted with 0 to 5 substituents independently chosen from (a), (b), and (c) where

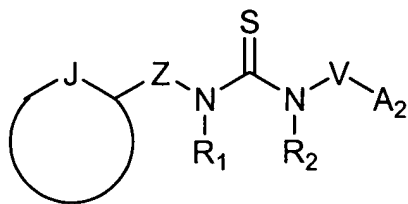
(a) is chosen from halogen, hydroxy, cyano, amino, nitro, oxo, $-\text{COOH}$, $-\text{CONH}_2$, $-\text{SO}_2\text{NH}_2$, $-\text{SH}$, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy,

(b) is chosen from C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_1 - C_6 alkoxy, C_2 - C_6 alkanoyl, and C_1 - C_8 alkoxycarbonyl, and

(c) is $-\text{GR}_a$ where G is chosen from $-(\text{CH}_2)_n-$, C_2 - C_4 alkenyl, C_2 - C_4 alkynyl, $-(\text{CH}_2)_n\text{O}(\text{CH}_2)_m-$, and $-(\text{CH}_2)_n\text{N}(\text{CH}_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and R_a is chosen from C_3 - C_8 cycloalkyl, piperidinyl, piperazinyl, morpholinyl, tetrahydroisoquinolinyl, indanyl, tetrahydronaphthyl, phenyl, pyridyl, benzothiophenyl, and benzofuranyl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(\text{C}_1$ - C_4 alkyl)amino, C_2 - C_4 alkanoyl, C_1 - C_4 alkoxycarbonyl, C_1 - C_2 haloalkyl, C_1 - C_2 haloalkoxy, and phenyl.

73. A compound of Formula II



Formula II

or a pharmaceutically acceptable salt thereof wherein

A₂ is C₃-C₈ cycloalkyl, a partially unsaturated or aromatic carbocyclic group, a saturated, partially unsaturated, or an aromatic heterocyclic group substituted with 0 to 5 substituents independently chosen from:

- (a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -CONH₂, -SO₂NH₂, -SH, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy, and
- (b) C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₁-C₆alkoxy, C₂-C₆alkenyloxy, C₁-C₄alkoxy(C₁-C₄alkyl), amino(C₁-C₆alkyl), mono- and di-(C₁-C₆alkyl)amino, mono- and di-(C₁-C₄alkyl)aminoC₁-C₄alkyl, C₂-C₆alkanoyl, C₂-C₈alkanoyloxy, C₁-C₈alkoxycarbonyl, -mono- and di-(C₁-C₆alkyl)carboxamide, (C₃-C₇cycloalkyl)carboxamide, mono- and di-(C₁-C₆alkyl)sulfonamide, C₁-C₆alkylthio, aryl(C₀-C₄alkyl)thio, C₁-C₆alkylsulfinyl, and C₁-C₆alkylsulfonyl, and

(c) -GR_a where

G is chosen from -(CH₂)_n-, C₂-C₄alkenyl, C₂-C₄alkynyl, -(CH₂)_nO(CH₂)_m-, and -(CH₂)_nN(CH₂)_m-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C₃-C₈cycloalkyl, C₂-C₇monocyclic heterocycloalkyl, C₅-

C₁₀bicyclic heterocycloalkyl, indanyl, tetrahydronaphthyl, aryl, and heteroaryl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from

halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₂-C₄alkanoyl, C₁-C₄alkoxycarbonyl, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, and phenyl;

V is C₁-C₆ alkyl, C₂-C₆alkenyl, or absent; and

Z is carbonyl, thiocarbonyl, or imino;

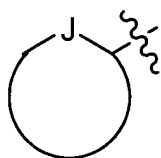
R₁ and R₂ are independently

hydrogen, or

C₁-C₆alkyl, C₂-C₆ alkenyl, or C₂-C₆ alkynyl, each of which is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkoxy, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy, or

R₁ and R₂ are joined to form a 5- to 7-membered saturated or mono-unsaturated ring optionally containing one additional heteroatom chosen from N, S, and O, which 5- to 7-membered saturated or mono-unsaturated ring is substituted with 0 to 3 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, and C₁-C₂haloalkoxy;

the group:



is a group of Formula (i)

that is a saturated, partially unsaturated, or aromatic heterocyclic group where J is O, S, or NR₃ substituted with 0 to 5 substituents independently chosen from: (a), (b), and (c) above; and

R₃ is

(d) hydrogen, C₁-C₂haloalkyl, or C₁-C₂haloalkoxy;

(e) C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₁-C₄alkoxy(C₁-C₄alkyl), or amino(C₁-C₆)alkyl, or

(f) -LR_b where

L is chosen from -(CH₂)_r-, C₂-C₄alkenyl, C₂-C₄alkynyl, -(CH₂)_rO(CH₂)_s-, and -(CH₂)_rN(CH₂)_s-, where r and s are independently 0, 1, 2, 3, or 4; and

R_b is chosen from C₃-C₈cycloalkyl, C₂-C₇monocyclic heterocycloalkyl, C₅-C₁₀bicyclic heterocycloalkyl, indanyl, tetrahydronaphthyl, aryl, and heteroaryl;

each of which (e) and (f) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, and phenyl.

74. A compound or salt according to Claim 73 wherein Z is carbonyl.

75. A compound or salt according to Claim 73 or Claim 74 wherein V is absent or V is C₁-C₄alkyl.

76. A compound or salt according to any one of Claims 73 to 75 wherein R_1 and R_2 are independently hydrogen or methyl.

77. A compound or salt according to any one of Claims 73 to 76 wherein A_2 is C_5 - C_7 cycloalkyl, phenyl, pyridyl, naphthyl, pyrimidinyl, pyrazinyl, benzothiazolyl, benzodioxyl, quinolinyl, or isoquinolinyl, each of which is substituted with 0 to 5 substituents independently chosen from (a), (b), and (c).

78. A compound or salt according to any one of Claims 73 to 77 wherein A_2 is C_5 - C_7 cycloalkyl, phenyl, pyridyl, naphthyl, benzothiazolyl, benzodioxyl, quinolinyl, or isoquinolinyl, each of which is substituted with 0 to 5 substituents independently chosen from

(a) halogen, hydroxy, cyano, amino, nitro, oxo, $-COOH$, $-CONH_2$, $-SO_2NH_2$, $-SH$, C_1 - C_2 haloalkyl, and C_1 - C_2 haloalkoxy,

(b) C_1 - C_6 alkyl, C_2 - C_6 alkenyl, C_2 - C_6 alkynyl, C_1 - C_6 alkoxy, C_2 - C_6 alkanoyl, and C_1 - C_8 alkoxycarbonyl, and

(c) $-GR_a$ where

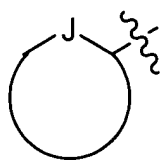
G is chosen from $-(CH_2)_n-$, C_2 - C_4 alkenyl, C_2 - C_4 alkynyl, $-(CH_2)_nO(CH_2)_m-$, and $-(CH_2)_nN(CH_2)_m-$, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C_3 - C_8 cycloalkyl, piperidinyl, piperazinyl, morpholinyl, tetrahydroisoquinolinyl, indanyl, tetrahydronaphthyl, phenyl, pyridyl, benzothiophenyl, and benzofuranyl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from

halogen, hydroxy, amino, C_1 - C_4 alkyl, C_1 - C_4 alkoxy, mono- and di- $(C_1$ - C_4 alkyl)amino, C_2 - C_4 alkanoyl, C_1 - C_4 alkoxycarbonyl, C_1 - C_2 haloalkyl, C_1 - C_2 haloalkoxy, and phenyl.

79. A compound or salt according to any one of Claims 73 to 78 wherein



is a group of Formula (i)

where Formula (i) is a heteroaryl group that is pyridyl, pyrimidinyl, thienyl, pyrrolyl, furanyl, pyrazolyl, imidazolyl, thiazolyl, triazolyl, thiadiazolyl, oxazolyl, isoxazolyl, benzofuranyl, benzothiazolyl, benzothiophenyl, benzoxadiazolyl, benzo[*d*]oxazolyl, dihydrobenzodioxynyl, indolyl, pyrazolopyrimidinyl, or thienylpyrazolyl oriented such that the heteroatom J is adjacent to the point of attachment of the group of Formula (i)

the group of Formula (i) is substituted with 0 to 5 substituents independently chosen from: (a), (b), and (c);

J is S, O, or NR₃; and

R₃ is

(d) hydrogen, C₁-C₂haloalkyl, or C₁-C₂haloalkoxy;

(e) C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₁-C₄alkoxy(C₁-C₄alkyl), or amino(C₁-C₆)alkyl, or

(f) -LR_b where

L is chosen from -(CH₂)_r-, C₂-C₄alkenyl, C₂-C₄alkynyl, -(CH₂)_rO(CH₂)_s-, and

-(CH₂)_rN(CH₂)_s-, where r and s are independently 0, 1, 2, 3, or 4; and

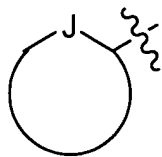
R_b is chosen from C₃-C₈cycloalkyl, C₂-C₇monocyclic heterocycloalkyl, C₅-

C₁₀bicyclic heterocycloalkyl, indanyl, tetrahydronaphthyl, aryl, and heteroaryl;

each of which (e) and (f) is substituted with 0 to 5 substituents independently chosen from

halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, and phenyl.

80. A compound or salt of Claim 79 wherein
wherein



is a group of Formula (i)

where Formula (i) is a heteroaryl group that is pyridyl, pyrimidinyl, thienyl, pyrrolyl, furanyl, pyrazolyl, imidazolyl, thiazolyl, triazolyl, thiadiazolyl, oxazolyl, isoxazolyl, benzofuranyl, benzothiazolyl, benzothiophenyl, benzoxadiazolyl, benzo[*d*]oxazolyl, dihydrobenzodioxynyl, indolyl, pyrazolopyrimidinyl, or thienylpyrazolyl oriented such that the heteroatom J is adjacent to the point of attachment of the group of Formula (i) in Formula II;

the group of Formula (i) is substituted with 0 to 5 substituents independently chosen from:

(a) halogen, hydroxy, cyano, amino, nitro, oxo, -COOH, -CONH₂, -SO₂NH₂, -SH, C₁-

C₂haloalkyl, and C₁-C₂haloalkoxy, and

(b) C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₁-C₆alkoxy, C₂-C₆alkanoyl, and C₁-C₈alkoxycarbonyl,

(c) -GR_a where

G is chosen from -(CH₂)_n-, C₂-C₄alkenyl, C₂-C₄alkynyl, -(CH₂)_nO(CH₂)_m-, and -(CH₂)_nN(CH₂)_m-, where n and m are independently 0, 1, 2, 3, or 4; and

R_a is chosen from C₃-C₈cycloalkyl, piperidinyl, piperazinyl, morpholinyl, tetrahydroisoquinolinyl, indanyl, tetrahydronaphthyl, phenyl, pyridyl, benzothiophenyl, and benzofuranyl;

each of which (b) and (c) is substituted with 0 to 5 substituents independently chosen from

halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₂-C₄alkanoyl, C₁-C₄alkoxycarbonyl, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, and phenyl;

J is S, O, or NR₃, and

R₃ is

(d) hydrogen,

(e) C₁-C₆alkyl, or

(f) -LR_b where

L is chosen from -(CH₂)_r-, -(CH₂)_rO(CH₂)_s-, and -(CH₂)_rN(CH₂)_s-, where r and s are independently 0, 1, 2, 3, or 4; and

R_b is chosen from C₃-C₈cycloalkyl, piperidinyl, piperazinyl, morpholinyl, indanyl, tetrahydronaphthyl, phenyl, and pyridyl;
 each of which (e) and (f) is substituted with 0 to 5 substituents independently chosen from halogen, hydroxy, amino, C₁-C₄alkyl, C₁-C₄alkoxy, mono- and di-(C₁-C₄alkyl)amino, C₁-C₂haloalkyl, C₁-C₂haloalkoxy, and phenyl.

81. A compound or pharmaceutically acceptable salt thereof, in which the compound is selected from

1-(Furan-2-carbonyl)-3-(4-benzo[*d*]thiazol-2-yl-phenyl)-thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-[5-(benzo[*d*]oxazol-2-yl)-2-methyl]phenylthiourea;
 1-(3-(Benzo[*d*]thiazol-2-yl)phenyl)-3-(2-phenoxyacetyl) thiourea;
 1-(4-(Benzo[*d*]oxazol-2-yl)phenyl)-3-propionylthiourea;
 1-(Pyridin-3-carbonyl)-3-(4-benzo[*d*]thiazol-2-yl-phenyl)-thiourea;
 1-[3-(2-chlorophenyl-5-methyl-isoxazol-4-yl)-carbonyl]-3-(4-isopropylphenyl)thiourea;
 Butyl 4-(3-(2-phenoxyacetyl) thioureido)benzoate;
 Butyl 4-(3-acetylthioureido)benzoate;
 Butyl 4-(3-(2-(3-chlorophenoxy) acetyl) thioureido)benzoate;
 Butyl 4-(3-(3-phenoxypropanoyl) thioureido)benzoate;
 Butyl 4-(3-(2-(naphthalen-3-yloxy)acetyl)thioureido)benzoate;
 Butyl 4-(3-(benzofuran-2-yl-carbonyl)thioureido)benzoate;
 Ethyl 2-(4-(3-(2-phenoxyacetyl)thioureido)phenyl)acetate;
 Ethyl 4-(3-(2-phenoxyacetyl)thioureido)benzoate;
 Butyl 4-(3-(2-methoxyacetyl) thioureido)benzoate;
 Butyl 4-(3-(2-(2,4-dichlorophenoxy) acetyl)thioureido)benzoate;
 Butyl 4-(3-(2-(4-tert-butylphenoxy) acetyl)thioureido)benzoate;
 Butyl 4-(3-(2-(4-(benzyloxy) phenoxy)acetyl)thioureido)benzoate;
 Butyl 4-(3-(2-(2-methoxyphenoxy) acetyl)thioureido)benzoate;
 Butyl 4-(3-(2-(*o*-tolylloxy)acetyl)thioureido)benzoate;
 Butyl 4-(3-(2-(2,4,6-trichlorophenoxy)acetyl)thioureido)benzoate;
 Butyl 4-(3-(3,4-dichlorophenyl) carbonyl)thioureido)benzoate;
 1-(3,4-dichlorophenyl-carbonyl)-3-(3-trifluoromethylphenyl)thiourea;

1-(3,4-Dichlorophenyl-carbonyl)-3-(3-benzoxo-phenyl)thiourea;
 1-(3,4-Dichlorophenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(3,4-Difluorophenyl-carbonyl)-3-(3-(5-methylfuran-2-yl)-phenyl)thiourea;
 Butyl 4-(3-(naphth-2-yl) carbonyl)thioureido)benzoate;
 Butyl 4-(3-(4-cyanophenyl) carbonyl)thioureido)benzoate;
 Butyl 4-(3-(methylacetate) carbonyl)thioureido)benzoate;
 1-((Benzofuran-2-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;
 Butyl 4-(3-(2-(3,4-dichlorophenoxy)acetyl)thioureido)benzoate;
 1-(4-Butylphenyl)-3-(2-phenoxyacetyl)thiourea;
 (Amino-(3-(benzyloxy)phenyl) methanethiocarbamoyl)methyl acetate;
 1-(3-(Methylthio)propanoyl)-3-(3-(benzyloxy)phenyl)thiourea;
 1-(2-(2-Chlorophenoxy)acetyl)-3-(3-(benzyloxy)phenyl)thiourea;
 Butyl 4-(3-(naphth-1-yl) carbonyl)thioureido)benzoate;
 (S)-1-(Amino-*N*-p-(butylacetate) methanethiocarbamoyl)ethyl acetate;
 Butyl 4-(3-(2-(2-methoxyethoxy)acetyl)thioureido)benzoate;
 (Amino-*N*-(4-cyclohexylphenyl)methanethiocarbamoyl)(phenyl)methyl acetate;
 Ethyl 3-(amino-*N*-(4-cyclohexylphenyl)methanethiocarbamoyl)propanoate;
 1-Butyryl-3-(4-cyclohexylphenyl)thiourea;
 (S)-1-(Amino-*N*-(4-cyclohexylphenyl)methanethiocarbamoyl)ethyl acetate;
 1-(3-(Benzyloxy)phenyl)-3-(2-hydroxyacetyl)thiourea;
 Butyl 4-(3-(2-(2,6-dichlorophenoxy) acetyl)thioureido)benzoate;
 Butyl 4-(3-(2-(3-methoxyphenoxy) acetyl) thioureido)benzoate;
 1-[(1-methylimidazol-2-yl) –carbonyl]-3-(3-benzoxo-phenyl)thiourea;
tert-Butyl 2-(amino-*N*-(3-(benzyloxy) phenyl)methanethiocarbamoyl) pyrrolidone-1-carboxylate;
 Butyl 4-(3-(pyrrolidin-1-yl) carbonyl)thioureido)benzoate;
 Butyl 4-(3-(1-methyl-benzofuran-2-yl) carbonyl)thioureido)benzoate;
 1-(4-Hexylphenyl)-3-(2-phenoxyacetyl)thiourea;
 1-(4-(Pentyloxy)phenyl)-3-(2-phenoxyacetyl)thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(4-pentyloxy)-phenyl)thiourea;
 1-(4-Pentylphenyl)-3-(2-phenoxyacetyl)thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(4-pentyl)-phenyl)thiourea;

1-((Benzofuran-2-yl-carbonyl)-3-(4-pentyloxy)-phenyl)thiourea;
 1-(4-Butoxyphenyl)-3-(2-phenoxyacetyl)thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(3-phenyl)-phenyl)thiourea;
 1-(2-Phenoxyacetyl)-3-(3-phenyl)-phenylthiourea;
 Isopropyl 4-(3-(benzofuran-2-yl) carbonyl)thioureido)benzoate;
 1-(2-phenoxyacetyl)-3-94-fluoro-phenyl)-phenylthiourea;
 1-(3-benzylphenyl)-3-(2-phenoxyacetyl)thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(3-benzyl)-phenyl)thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(4-benzyl)-phenyl)thiourea;
 1-(4-(p-Tolyloxy)phenyl)-3-(2-phenoxyacetyl)thiourea;
 Isobutyl 4-(3-(benzofuran-2-yl) carbonyl)thioureido)benzoate;
 Isobutyl 4-(3-(2-phenoxyacetyl)thioureido)benzoate;
 1-(2-(phenylmethanone)phenyl)-3-(2-phenoxyacetyl)thiourea;
 1-(3-(Phenylcarbamoyl)phenyl)-3-(2-phenoxyacetyl)thiourea;
 1-(3-(2-Methylpyrimidin-4-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;
 1-(4-(4-Chlorophenoxy)phenyl)-3-(2-phenoxyacetyl)thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(4-Chlorophenoxy)-phenyl)thiourea;
 1-(4-(3,4-Dihydroisoquinolin-2(1H)-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;
 1-(3-Fluoro-4-(octahydroquinolin-1(2H)-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;
 1-(3-Fluoro-4-(octahydroquinolin-1(2H)-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;
 1-(3-Fluoro-4-(piperidin-1-yl)phenyl)-3-(2-phenoxyacetyl)thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(3-fluoro-4-(piperidin-1-yl)phenyl)-phenyl)thiourea;
 1-(3-(3-Methoxybenzyloxy) phenyl)-3-(2-phenoxyacetyl) thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(3-(3-methoxybenzyloxy)) phenyl)-phenyl)thiourea;
 1-(3-(2-Methoxybenzyloxy) phenyl)-3-(2-phenoxyacetyl) thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(3-(2-Methoxybenzyloxy) phenyl)-thiourea;
 1-(3-(4-Methoxybenzyloxy) phenyl)-3-(2-phenoxyacetyl) thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(3-(cyclohexylmethoxy) phenyl)-thiourea;
 1-(3-(Cyclohexylmethoxy) phenyl)-3-(2-phenoxyacetyl)thiourea;
 1-((Benzofuran-2-yl-carbonyl)-3-(4-(5,6-dihydropyridin-1(2H)-yl))phenyl)-thiourea;
 1-((5-Methoxy-benzofuran-2-yl-carbonyl)-3-(3-benzyloxy-phenyl)-thiourea;

Butyl 4-(3-(5-chloro-benzofuran-2-yl) carbonyl)thioureido)benzoate;
 1-(7-Methoxy-benzofuran-2-yl-carbonyl)-3-(3-methoxybenzyloxy)phenyl)-phenyl)thiourea;
 1-(2,3,4-tetrahydronaphthalen-2-yl-carbonyl)-3-(3-(3-methoxybenzyloxy))phenyl)-
 phenyl)thiourea;
 1-(2-(4-(Trifluoromethoxy) phenoxy) acetyl)-3-(3-(benzyloxy)phenyl) thiourea;
 1-(3-(Benzyloxy)phenyl)-3-(2-(pyridin-3-yloxy)acetyl) thiourea;
 1-(4-oxo-4-*H*-chromen-2-yl-carbonyl)-3-(3-methoxybenzyloxy)phenyl)-phenyl)thiourea;
 1-(3-(Benzyloxy)phenyl)-3-(2-(pyridin-2-yloxy)acetyl) thiourea;
 1-(Pyridin-2-yl-carbonyl)-3-(3-methoxybenzyloxy)phenyl)-phenyl)thiourea;
 1-(3-Chloro-benzo[*b*]thiophen-2-yl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;
 1-(4-Trifluoromethoxy-phenyl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;
 1-(5-Methylisoxazol-3-yl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;
 1-(2-Methyl-5-phenyl-furan-3-yl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;
 1-(4-Trifluoromethyl-phenyl-carbonyl)-3-(3-ethoxybenzyloxy)phenyl)-phenyl)thiourea;
 1-(3-Chloro-benzo[*b*]thiophen-2-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;
 1-(4-Trifluoromethoxyphenyl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;
 1-(3,5-Dimethylisoxazol-4-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;
 1-(5-Methylisoxazol-3-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;
 1-(4-Trifluoromethylphenyl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;
 (S)-1-(Amino-*N*-(3-(benzyloxy) phenyl)methanethiocarbamoyl)ethyl acetate;
 (S)-1-(Amino-*N*-(3-phenoxyphenyl)methane thiocarbamoyl)ethyl acetate;
 (S)-1-(Amino-*N*-(3-benzyl-phenyl)methanethiocarbamoyl)ethyl acetate;
 Ethyl 1-(2-fluoro-4-(3-(benzofuran-2-yl-carbonyl)thioureido)phenyl)-4-phenylpiperidine-4-
 carboxylate;
 1-(4-Trifluoromethylphenyl-carbonyl)-3-(2-phenylbenzo[*d*][1,3]dioxol-6-yl) phenyl)thiourea;
 1-(3-Chloro-methylbenzo[*b*]thiophen-2-yl-carbonyl)-3-(2-phenylbenzo[*d*][1,3]dioxol-6-yl)
 phenyl)thiourea;
 1-(4-Trifluoromethoxyphenyl-carbonyl)-3(2-phenylbenzo[*d*][1,3]dioxol-6-yl) thiourea;
 1-(5-Methylisoxazol-3-yl-carbonyl)-3-(2-phenylbenzo[*d*][1,3]dioxol-6-yl) phenyl)thiourea;
 1-(3-((*R*)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)thiourea;
 1-(3-((*R*)-1-Phenylethoxy)phenyl)-3-(3-chloro-methylbenzo[*b*]thiophen-2-yl-carbonyl)thiourea;

1-(3-((R)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethoxy-phenyl-carbonyl)thiourea;
 1-(3-((R)-1-Phenylethoxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)thiourea;
 1-(3-((S)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)-thiourea;
 1-(3-((S)-1-Phenylethoxy)phenyl)-3-(4-trifluoromethoxyphenyl-carbonyl)-thiourea;
 1-(3-((S)-1-Phenylethoxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)-thiourea;
 1-(3-(Phenethyloxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)-thiourea;
 1-(3-(Phenethyloxy)phenyl)-3-(3-chloro-methylbenzo[b]thiophen-2-yl-carbonyl)-thiourea;
 1-(3-(Phenethyloxy)phenyl)-3-(4-trifluoromethoxyphenyl-carbonyl)-thiourea;
 1-(3-(Phenethyloxy)phenyl)-3-(3,5-dimethylisoxazol-4-yl-carbonyl)-thiourea;
 1-(3-(Phenethyloxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)-thiourea;
 1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(4-trifluoromethylphenyl-carbonyl)thiourea;
 1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(4-trifluoromethoxyphenyl-carbonyl)thiourea;
 1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(3,5-dimethylisoxazol-4-yl-carbonyl)thiourea;
 1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(5-methylisoxazol-3-yl-carbonyl)thiourea;
 1-(2-Phenylbenzo[d][1,3]dioxol-6-yl)-3-(benzofuran-2-yl-carbonyl)thiourea;
 1-(3-((S)-1-Phenylethoxy)phenyl)-3-(benzofuran-2-yl-carbonyl)-thiourea;
 1-(3-(Phenethyloxy)phenyl)-3-(benzofuran-2-yl-carbonyl)-thiourea;
 1-(3-(2,3-Dihydro-1H-inden-2-yloxy)phenyl)-3-(benzofuran-2-yl-carbonyl)thiourea;
 1-(3-((R)-1-Phenylethoxy)phenyl)-3-(benzofuran-2-yl-carbonyl)thiourea;
 1-(2,4-dimethylthiazol-5-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;
 1-(1-methyl-pyrrol-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;
 1-(3-(Phenethyloxy)phenyl)-3-((2,7-dimethylpyrazolo[1,5-a]pyrimidin-6-yl)-carbonyl)-thiourea;
 1-(3-(Phenethyloxy)phenyl)-3-(1-ethyl-3-methyl-1*H*-pyrazol-5-yl-carbonyl)-thiourea;
 1-(3-(2,3-Dihydro-1*H*-inden-2-yloxy)phenyl)-3-(2,4-dimethylthiazol-5-yl-carbonyl)thiourea;
 1-(3-(2,3-Dihydro-1*H*-inden-2-yloxy)phenyl)-3-((2,7-dimethylpyrazolo[1,5-a]pyrimidin-6-yl)-carbonyl)thiourea;
 1-(3-(2,3-Dihydro-1*H*-inden-2-yloxy)phenyl)-3-(1-ethyl-3-methyl-1*H*-pyrazol-5-yl-carbonyl)thiourea;
 1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;
 1-(Benzo[d]thiazol-2-yl-carbonyl)-3-(3-phenyloxy)phenyl)-phenyl)thiourea;
 1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;

1-(Benzo[*d*]thiazol-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(3-fluoro-5-pentoxy-phenyl)thiourea;
 1-(2-Methyl-pyridin-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-(1-phenyl-1*H*-pyrazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-(1-phenyl-1*H*-pyrazol-5-yl)-3-((3-phenyloxy)phenyl)-phenylthiourea;
 1-(2-Phenylbenzo[*d*][1,3]dioxol-6-yl)-3-(1-phenyl-1*H*-pyrazol-5-yl-carbonyl)thiourea;
 1-(1-phenyl-1*H*-pyrazol-5-yl)-3-(4-pentoxy-phenyl)-phenylthiourea;
 1-(1-phenyl-1*H*-pyrazol-5-yl)-3-((3-phenyloxy-phenyl)-phenyl)thiourea;
 1-(Methylbenzo[*b*]thiophen-2-yl-carbonyl)-3-((4-pentoxy-phenyl)-phenyl)thiourea;
 1-(isoxazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-(isoxazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;
 1-(isoxazol-5-yl-carbonyl)-3-(4-(pentyl) phenyl)thiourea;
 1-(isoxazol-5-yl-carbonyl)-3-(4-(pentoxy) phenyl)thiourea;
 1-(isoxazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;
 1-(isoxazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;
 1-((3-(trifluoromethyl)phenyl)furan-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-((3-(trifluoromethyl)phenyl)furan-2-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;
 1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;
 1-(5-Nitro-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-(5-Nitro-benzofuran-2-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;
 1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(4-pentyl phenyl)thiourea;
 1-(5-Bromo-benzofuran-2-yl-carbonyl)-3-(4-pentoxy phenyl)thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(4-(hex-1-ynyl)phenyl) thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(3-(2-(pyridin-3-yl)ethynyl)phenyl) thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(3-(2-(pyridin-3-yl)ethyl)phenyl) thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(3-(pyridin-2-yl) phenyl) thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(3-(pyridin-2-yl-oxy)-phenyl) thiourea;
 1-(3,5-Dimethylisoxazole-4-yl-carbonyl)-3-(4-(hex-1-ynyl)phenyl) thiourea;
 1-(3,5-Dimethylisoxazole-4-yl-carbonyl)-3-(3-(2-(pyridin-3-yl)ethynyl)phenyl) thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-((3-(trifluoromethyl)benzyloxy)-phenyl) thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(4-((1-methylpiperidin-4-yl)methoxy)-3-fluorophenyl)thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(3-(trifluoromethyl)-4-(piperidin-1-yl)phenyl)thiourea;
 1-(1,3-dimethyl-1*H*-thieno[2,3-*c*]pyrazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;
 1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(5-(2-methylthiazol-4-yl)isoxazol-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;
 1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;
 1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(1,5-Dimethyl-1*H*-pyrazol-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;
 1-(1-Methyl-3-*tert*-butyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-(1-Methyl-3-*tert*-butyl-1*H*-pyrazol-3-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;
 1-(1-Methyl-3-*tert*-butyl-1*H*-pyrazol-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(2-Trifluoromethyl-5-methyl-furan-2-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(2-Trifluoromethyl-5-methyl-furan-2-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;
 1-(2-Trifluoromethyl-5-methyl-furan-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(3-(benzyloxy) phenyl)thiourea;
 1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(3-(phenoxy) phenyl)thiourea;
 1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(4-(pentyl) phenyl)thiourea;
 1-(Benzo[*c*][1,2,5]oxadiazol-5-yl-carbonyl)-3-(4-(pentoxy) phenyl)thiourea;
 1-(2,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(2,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;
 1-(2,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-6-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;
 1-(3-Methylisoxazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(3-Methylisoxazol-4-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;
 1-(3-Methylisoxazol-4-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;
 1-(3-Methylisoxazol-4-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;
 1-(5-Methyl-2-phenyl-2*H*-1,2,3-triazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(5-Methyl-2-phenyl-2*H*-1,2,3-triazol-4-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;
 1-(5-Chloro-3-methylbenzo[*b*]thiophen-2-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;

1-(5-Chloro-3-methylbenzo[*b*]thiophen-2-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;
 1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;
 1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;
 1-(1,3-Dimethyl-1*H*-pyrazol-5-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;
 1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(3-(phenyloxy)-phenyl)thiourea;
 1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(4-(pentyl)-phenyl)thiourea;
 1-(2-(Pyridin-3-yl)thiazol-4-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;
 1-(4-Methoxy-benzofuran-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;
 1-(4-Methoxy-benzofuran-2-yl-carbonyl)-3-(3-fluoro-4-pentoxy-phenyl)thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(3,5-dibromo-4-(pent-4-enyloxy)phenyl)thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(3-((pyridin-3-yl)methyl) phenyl)thiourea;
 1-(5-Iodo-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;
 1-(5-Phenyl-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;
 1-(5-(2-Pyridyl)benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)phenyl)thiourea;
 1-(3-Propoxy-pyridin-2-yl-carbonyl)-3-(4-(pentyl)phenyl)thiourea;
 1-(2,5-Dichlorothiophen-3-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;
 1-(3-Methyl-5-(methylthio)-4-vinylthiophen-2-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;
 1-(3-Methyl-5-(methylthio)-4-vinylthiophen-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;
 1-(5-(Methylthio)-thiophen-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;
 1-(5,7-Dimethylpyrazolo[1,5-*a*]pyrimidin-2-yl-carbonyl)-3-(4-(benzyloxy)-phenyl)thiourea;
 1-(2,5-Dichlorothiophen-3-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;
 1-(3-Methyl-5-(methylthio)-4-vinylthiophen-2-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;
 1-(5-(Methylthio)-thiophen-2-yl-carbonyl)-3-(4-(phenoxy)-phenyl)thiourea;
 1-(7-Fluoro-benzofuran-2-yl-carbonyl)-3-(4-(pentoxy)-phenyl)thiourea;
 1-(7-Fluoro-benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(2-(1,3-Dioxoisindolin-2-yl)acetyl)-3-(3-phenoxyphenyl)thiourea;
 1-(2-(1,3-Dioxoisindolin-2-yl)acetyl)-3-(3-benzyloxyphenyl)thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(3-(benzyloxy)methyl-phenyl)thiourea ;
 1-(Benzofuran-2-yl-carbonyl)-3-(3-(phenylamino)methyl-phenyl)thiourea;

1-(Benzofuran-2-yl-carbonyl)-3-(4-(*N*-benzyl-*N*-methylamino)-3-fluorophenyl) thiourea;
 Phenyl 3-(3-((benzofuran-2-yl)-carbonyl)thioureido)benzoate;
 1-(4-Cyanophenyl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(4-Cyanophenyl-carbonyl)-3-(4-(pentyloxy)-phenyl)thiourea;
 1-(4-Cyanophenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(Quinoxalin-2-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(Quinoxalin-2-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;
 1-(Quinoxalin-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(4-Trifluoromethoxyphenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(4-Trifluoromethoxyphenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;
 1-(4-Trifluoromethylphenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;
 1-(4-Trifluoromethylphenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(5-Cyano-benzofuran-2-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(5-Cyano-benzofuran-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;
 1-(4-Methyl-1,2,3-thiadiazol-5-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;
 Butyl 4-(3-(4-Methyl-1,2,3-thiadiazol-5-yl-) carbonyl)thioureido) benzoate;
 1-(Pyrazin-2-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(Pyrazin-2-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(Pyrazin-2-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;
 1-(3-Fluoro-4-(pentyloxy)phenyl)-3-(2-(1,3-dioxoisindolin-2-yl)acetyl)thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(3-trifluoromethyl-4-pentoxy-phenyl)thiourea;
 1-(1-Benzyl-1*H*-tetrazol-5-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(4-(*N*-methyl-*N*-pentylamino)-3-fluorophenyl)thiourea;
 1-(1-Benzyl-1*H*-tetrazol-5-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(6-Trifluoromethyl-pyrid-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(6-Trifluoromethyl-pyrid-3-yl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;
 1-(3-Trifluoromethyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(3-Trifluoromethyl-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(3-Trifluoromethyl-phenyl-carbonyl)-3-(4-pentoxy-phenyl)thiourea;
 1-(3-Trifluoromethoxy-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;

1-(3-Trifluoromethoxy-phenyl-carbonyl)-3-(4-pentoxo-phenyl)thiourea;
 1-(2-Chloro-5-trifluoromethoxy-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(3-Difluoromethyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(3-Difluoromethyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(5-(Trifluoromethyl)-2-phenyloxazol-4-yl-carbonyl)-3-(3-(benzyloxy)-phenyl)thiourea;
 1-(5-(Trifluoromethyl)-2-phenyloxazol-4-yl-carbonyl)-3-(3-(phenoxy)-phenyl)thiourea;
 1-(5-(2-Chloro-5-trifluoromethylphenyl)-furan-2-yl-carbonyl)-3-(4-(pentoxo)-phenyl)thiourea;
 1-(3-Trifluoromethyl-4-methoxy-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(3-Trifluoromethyl-4-chloro-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(3-Trifluoromethyl-4-chloro-phenyl-carbonyl)-3-(4-pentoxo-phenyl)thiourea;
 1-(3-Trifluoromethyl-4-methyl-phenyl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea;
 1-(3-Trifluoromethyl-4-methyl-phenyl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(3-Trifluoromethyl-4-methyl-phenyl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-((5-Acetamidobenzofuran-2-yl)carbonyl)-3-(3-phenoxyphenyl)thiourea;
 1-Acetyl-3-(3-phenoxyphenyl)thiourea;
 1-Acetyl-3-(4-(pentylloxy)phenyl)thiourea;
 1-Acetyl-3-(4-pentylphenyl)thiourea;
 1-(Dimethylamino-acetyl)-3-(3-phenoxyphenyl)thiourea;
 1-(Dimethylamino-acetyl)-3-(3-benzyloxyphenyl)thiourea;
 1-(3,5-Dimethylisoxazole-4-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-((2,3,4,5,6-penta-fluorophenoxy)-phenyl)thiourea;
 1-(Benzofuran-2-yl-carbonyl)-3-(9-methyl-9*H*-fluoren-7-yl)thiourea;
 Pentyl 2-phenyl- 4-(3-(benzofuran-2-yl)thioureido)benzoate;
 1-(3-Pyrid-3-yl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea;
 1-(3-Pyrid-3-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea;
 1-(3-Pyrid-3-yl-carbonyl)-3-(4-pentoxo-phenyl)thiourea;
 1-(3-Pyrid-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea;
 1-(4-Phenylbutanoyl)-3-(3-phenoxyphenyl)thiourea;
 1-(4-Phenylbutanoyl)-3-(3-benzyloxyphenyl)thiourea;
 1-(2-Morpholinoacetyl)-3-(3-phenoxyphenyl)thiourea;
 1-(2-Morpholinoacetyl)-3-(4-(pentylloxy)phenyl)thiourea;

1-(2-Morpholinoacetyl)-3-(4-(pentyl)phenyl)thiourea;
1-(4-(Pentyloxy)phenyl)-3-(2-(piperidin-1-yl)acetyl)thiourea;
1-(*N*-Methyl-*N*-phenylamino-acetyl)-3-(3-benzyloxyphenyl)thiourea;
1-(Benzofuran-2-yl-carbonyl)-3-(6-pentoxypyrid-3-yl)thiourea;
1-(3-Pyrid-3-yl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea hydrochloride;
1-(3-Pyrid-3-yl-carbonyl)-3-(3-phenoxy-phenyl)thiourea hydrochloride;
1-(3-Pyrid-3-yl-carbonyl)-3-(4-pentyl-phenyl)thiourea hydrochloride;
1-(Benzofuran-2-yl-carbonyl)-3-(trifluoromethylthio-phenyl)thiourea;
1-(3-(Piperidin-1-yl)propanoyl)-3-(4-pentylphenyl)thiourea;
1-(3-(Piperidin-1-yl)propanoyl)-3-(4-(pentyloxy)phenyl)thiourea;
1-(3-(Piperidin-1-yl)propanoyl)-3-(3-phenoxyphenyl)thiourea;
1-(3-Morpholinopropanoyl)-3-(4-(pentyloxy)phenyl)thiourea;
1-(1-Methylpiperidin-3-yl-carbonyl)-3-(4-(pentyloxy)phenyl)thiourea;
1-(1-Methylpiperidin-3-yl-carbonyl)-3-(4-(pentyloxy)phenyl)thiourea;
1-(2-(2-methylpiperidin-1-yl)acetyl)-3-(4-(pentyloxy)phenyl)thiourea;
1-(2-Oxo-4-phenyl-pyrrolidin-1-ylcarbonyl)-3-(3-benzyloxy-phenyl)thiourea; and
1-(5-Trifluoromethoxy-benzofuran-2-yl-carbonyl)-3-(3-benzyloxy-phenyl)thiourea.

82. A pharmaceutical composition comprising a compound or salt according to any one of Claims 1 to 81 together with a pharmaceutically acceptable carrier, diluent, or excipient.

83. A pharmaceutical composition according to Claim 82, wherein the composition is formulated as an injectable fluid, an aerosol, a cream, a gel, a pill, a capsule, a syrup, a tablet, ophthalmic solution, or a transdermal patch.

84. A package comprising a pharmaceutical composition of Claim 82 in a container and further comprising instructions for using the composition to treat a patient suffering from Hepatitis C infection.

85. A compound or salt according to Claim 1 that exhibits an EC₅₀ of less than 10 micromolar in a replicon assay of HCV replication.

86. A compound or salt according to Claim 1 that exhibits an EC₅₀ of less than 1 micromolar in a replicon assay of HCV replication.

87. A method for treating Hepatitis C infection comprising administering to a patient in need of such treatment a therapeutically effective amount of a compound or salt according to Claim 1.

88. The method of Claim 87 wherein the patient is a human patient.

89. The method of Claim 88 wherein the therapeutically effective amount is an amount sufficient to significantly decrease the number of HCV antibodies in the patient's blood or serum.

90. A method of inhibiting HCV replication *in vivo* comprising administering to a patient infected with HCV a concentration of a compound or salt according to Claim 1 sufficient to inhibit HCV replicon replication *in vitro*.